



UNITED STATES ENVIRONMENTAL PROTECTION AGENCY  
REGION 5 CENTRAL REGIONAL LABORATORY  
536 SOUTH CLARK STREET  
CHICAGO, ILLINOIS 60605

200002

EPA Region 5 Records Ctr.



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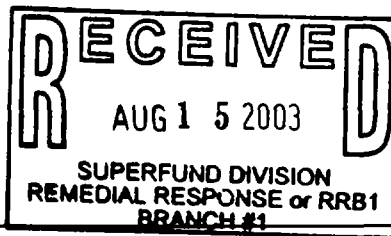
Print Date: 12/5/02

Subject: Review of Region 5 Data for Himco Dump

From: Xuyen Nguyen, Chemist *YN 12/5/02*  
Contractor to Region 5 Central Regional Laboratory  
Submitted to CRL on

To: Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago, IL 60604

Attached are Results for: Himco Dump



Analyses included in this report:

SVOA Standard List

*Sylvia Griffin*  
Data Management Coordinator and Date Received

12/05/02

Date Transmitted: 12/05/02

Please have the U.S. EPA Project Manager/Officer call the CRL Sample Coordinator at 3-7444 for any comments or questions.

Please sign and date this form below and return it with any comments to:

Sylvia Griffin  
Data Management Coordinator  
Region 5 Central Regional Laboratory  
ML-10C

\_\_\_\_\_/\_\_\_\_\_/\_\_\_\_\_  
Received by and Date

Comments:

SVDA

## **CASE NARRATIVE**

DATE: November 25, 2002

TDF#:05-0-047

Job#:249-0-1476-109-011-002

PROJECT NAME: Data set W/O # E2K0101 water samples for Semi-Volatile  
(CRL SOP GCMS026).

SITE NAME: HIMCO DUMP

ANALYST: Xuyen Nguyen, IITRI/ESAT

### **I CASE DESCRIPTION:**

ESAT received four water samples (E2K0101-02# 2003SY01R02, E2K0101-04# 2003SY01S04, E2K0101-05# 2003SY01D04 and E2K0101-06# 2003SY01S05) for semi-volatile analysis on November 01, 2002. The samples were extracted on November 4, 2002 using the continuous liquid/liquid extraction method per CRL SOP GCMS026. All samples extracts were analyzed on November 13, 2002. Their IS responses were below the lower QC limits. All The samples were re- analyzed on November 20, 2002 within the required hold time after discussing with the TOPO.

### **II. INSTRUMENT QUALITY CONTROLS:**

#### **1. Instrument Performance Checks:**

A GC/MS instrument performance check using DFTPP was made each day of analysis to determine if acceptable EPA tuning criteria were met. The QC criteria are the same as those found in CRL SOP GCMS026. All standards and samples were analyzed on GCMS #2. All QC criteria were met and no problems were observed.

#### **2. Initial Calibrations (IC):**

An acceptable five point IC is required for all target compounds before samples can be analyzed. The QC criterion for the IC states each analyte's %RSD must be < 30% for water (according to the criteria listed in CRL SOP GCMS026). All analytes met the requirements for the IC analyzed on November 20, 2002 except Benzoic acid (74.50%RSD), Hexachlorocyclopentadiene (53.7%RSD), 2,4-Dinitrophenol (58.6%RSD), 4,6-Dinitro-2-methylphenol(64.3%RSD) and Pentachlorophenol (94.8%RSD). These analytes were flagged estimated "J" for positive results and "UJ" for non-detects. The analyte 2,4-Dinitrophenol has three point calibration curve, 4-Nitrophenol has four point calibration curve and Pentachlorophenol has three point calibration curve. These compounds have an elevated detection limit.

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### **3. Continuing calibration (CC):**

An acceptable CC is required for all target compounds before samples can be analyzed. The QC criteria for the CC states each analyte's %D must be <25%. The continuing calibration, analyzed on November 20, 2002 reported all compounds within the quality control limits except 2,4-Dinitrophenol(52.4%D) and Pentachlorophenol (85.7%D). These analytes were flagged "J" for positive results and "UJ" for non-detects in the associated samples.

### **4. Internal Standard (IS) Area and Retention Time Summary:**

The QC criteria states that the IS area of the samples must range from 50 to 200% of the internal standard area of the corresponding CC. The RT of the IS for samples must also be within  $\pm 30$  seconds of the RT of the IS for the corresponding standard. All internal standards area met the QC requirements .

## **III. METHOD QUALITY CONTROL:**

### **1. Method Blank Results:**

A lab blank was prepared for each day samples were extracted to check the GC/MS, extraction glassware and the reagents for laboratory contamination.

The extraction blank EK21101-blk1 reported bis(2-Ethylhexyl)phthalate at 6ug/L . This analyte was qualified as non-detected "UJ" when the sample result is less than(<) ten (10) times the blank results and detected "J" when the sample results at or above the reporting limit(RL) in all the associated samples.

### **2. Surrogate Spike Compound Results:**

All surrogate recoveries were within the control limits stated in the CRL SOP GCMS026 except in the sample E2I0501-MSD1, only one surrogate Nitrobenzene-d5(34%) was outside the control QC limits, no action was required.

### **3. Laboratory Control Sample (LCS) Results:**

The EK21101-BS1 and EK21101-BSD1 were each spiked with 500ul of the 100ug/ml ABN LCS standard solution. This corresponded to a 50ug spiked into one liter of water. In the EK21101-BS1 and EK21101-BSD1 samples, all the relative percent differences (RPD) were within QC limits except Dimethylphthalate(103%) and Diethylphthalate(52%). In the EK21101-BS1/BSD1 samples , all the percent recoveries were within limits except analytes Diethylphthalate(20%) and Pentachlorophenol(5%) which were outside limits in the BS1 sample but were within limits in the BSD1 sample so no action was required.. The analyte 2,4-Dinitrophenol was not detected in either the BS1 or BSD1. This analyte was flagged "Q" for detected compound and "R" for non-detected compound in all the samples.

### **4. Matrix Spike/Matrix Spike Duplicate (MS/MSD) Results**

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Sample E2K0101-06 was used as the MS1/MSD1 samples. The MS1 and MSD1 samples were each spiked with 500ul of the 100ug/ml ABN LCS standard solution.

In the MS1 sample all the percent recoveries were within limits except Pentachlorophenol (188%) and 3,3-Dichlorobenzidine(26%).

The analyte 2,4-Dinitrophenol was not detected in either MS1 or MSD1. This analyte was flagged "Q" for detected and "R" for non-detected compounds in the native sample.

In the MSD1 sample all the percent recoveries were within limits except: bis(2-Chloroethyl)ether (10%), 1,3-Dichlorobenzene(11%), 1,4-Dichlorobenzene(12%), 1,2-Dichlorobenzene(16%), bis(2-chloroisopropyl)ether(24%), 2-Nitrophenol(38%) N-Nitroso-di-n-propylamine(32%), Nitrobenzene(32%), Isophorone(42%), bis(2-Chloroethoxy)methane(44%), 1,2,4-Trichlorobenzene(32%) 4-Chloroaniline (26%), Pentachlorophenol(162%) and 3,3-Dichlorobenzidine (18%). All these compounds percent recoveries were within limits in the MS1 except the analytes Pentachlorophenol and 3,3-Dichlorobenzidine were outside limits in both MS1 and MSD1. These compounds were flagged "J" for detected and "UJ" for non-detected compounds in the native sample.

The relative percent differences(RPD) were within limits except bis(2-Chloroethyl)ether (117%), 1,3-Dichlorobenzene(123%), 1,4-Dichlorobenzene(120%), 1,2-Dichlorobenzene (103%), bis(2-chloroisopropyl)ether(83%), Hexachloroethane(97%), N-Nitroso-di-n-propylamine(64%), Nitrobenzene(58%), Isophorone(42%), 2,4-Nitrophenol (51%),4-Chloroaniline (76%) and bis(2-Ethylhexyl)phthalate(29%). No other problems were observed.

#### **IV.SAMPLES RESULTS:**

In the sample E2k0101-02 contained bis(2-Ethylhexyl)phthalate(5.0ug/L)

In the sample E2k0101-04 contained bis(2-Ethylhexyl)phthalate(7.0ug/L)

In the sample E2k0101-05 contained bis(2-Ethylhexyl)phthalate(4.0ug/L)

In the sample E2k0101-06 contained bis(2-Ethylhexyl)phthalate(6.0ug/L)

The raw /processed data files, method files and sequence files are archived on R5CRL\VOL2\XNGUYEN\2002.

Tentatively Identified Compounds (TICs) have been listed for each sample. These compounds listed on the Enviroforms were selected based on the best Mass spectrum match of the unknown compound to compounds in the library. These compounds are suggested and not confirmed. LIMS reports will be generated as soon as possible.

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Environmental Protection Agency Region 5  
**Central Regional Laboratory**

536 South Clark Street, Chicago, IL 60605  
Phone: (312) 353-8370 Fax: (312) 886-2591

**WORK ORDER**

Printed: 11/25/02 12:26:56P

**E2K0101**

**IITRI - ESAT Contract**

**Client:** Superfund, US EPA Region 5  
**Project:** Himco Dump

**Project Manager:** Jennifer Mokus  
**Project Number:** 2003SY01

**Report To:**

Howard Pham  
Superfund, US EPA Region 5

77 West Jackson Boulevard  
Chicago, IL 60604

Phone: (312) 353-2310  
Fax: (312) 886-6171

**Date Due:** Dec-03-02 15:00 (30 day TAT)

**Received By:** William Sargent

**Date Received:** Nov-01-02 11:27

**Logged In By:** William Sargent

**Date Logged In:** Nov-01-02 12:16

**Samples Received at:** °C  
**All containers intact:** No  
**Sample labels/COC agree:** No  
**Samples Preserved Properly:** No  
**Custody Seals Present:** No

Analysis	Due	TAT	Expires	Comments
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**E2K0101-01 2003SY01S01-SINK [Water] Sampled Oct-31-02 07:40 Central**

Volatiles Full List	Dec-03-02 12:00	30	Nov-07-02 07:40	
THF & 1,4-Dioxane	Dec-03-02 12:00	30	Nov-14-02 07:40	

**E2K0101-02 2003SY01R02-Pump Blank [Water] Sampled Oct-31-02 11:50 Central**

GFAA 5100 Cd	Dec-03-02 12:00	30	Apr-29-03 11:50	
Volatiles Full List	Dec-03-02 12:00	30	Nov-07-02 11:50	
SVOA Standard List	Dec-03-02 12:00	30	Nov-07-02 11:50	
Cyanide, Total	Dec-03-02 12:00	30	Nov-14-02 11:50	pH10
Hg Total CVAA	Dec-03-02 12:00	30	Nov-28-02 11:50	pH2
Solids, Dry Weight	Dec-03-02 12:00	30	Nov-07-02 11:50	For SVOA Standard List in batch EK21101
B ICP (W)	Dec-03-02 12:00	30	Apr-29-03 11:50	pH2
THF & 1,4-Dioxane	Dec-03-02 12:00	30	Nov-14-02 11:50	
Metals SF ICP (W)	Dec-03-02 12:00	30	Apr-29-03 11:50	pH2
GFAA SIMAA Metals	Dec-03-02 12:00	30	Apr-29-03 11:50	pH2

**E2K0101-03 2003SY01R03-Trip Blank [Water] Sampled Oct-31-02 11:50 Central**

Volatiles Full List	Dec-03-02 12:00	30	Nov-07-02 11:50	
THF & 1,4-Dioxane	Dec-03-02 12:00	30	Nov-14-02 11:50	

E2K0101

## IITRI - ESAT Contract

Client: Superfund, US EPA Region 5  
Project: Himco Dump

Project Manager: Jennifer Mokos  
Project Number: 2003SY01

Analysis	Due	TAT	Expires	Comments
<b>E2K0101-04 2003SY01S04-WT116A [Water] Sampled Oct-31-02 13:52 Central</b>				
GFAA 5100 Cd	Dec-03-02 12:00	30	Apr-29-03 13:52	
Volatiles Full List	Dec-03-02 12:00	30	Nov-07-02 13:52	
SVOA Standard List	Dec-03-02 12:00	30	Nov-07-02 13:52	
Cyanide, Total	Dec-03-02 12:00	30	Nov-14-02 13:52	pH10
Hg Total CVAA	Dec-03-02 12:00	30	Nov-28-02 13:52	pH2
Solids, Dry Weight	Dec-03-02 12:00	30	Nov-07-02 13:52	For SVOA Standard List in batch EK21101
B ICP (W)	Dec-03-02 12:00	30	Apr-29-03 13:52	pH2
THF & 1,4-Dioxane	Dec-03-02 12:00	30	Nov-14-02 13:52	
Metals SF ICP (W)	Dec-03-02 12:00	30	Apr-29-03 13:52	pH2
GFAA SIMAA Metals	Dec-03-02 12:00	30	Apr-29-03 13:52	pH2
<b>E2K0101-05 2003SY01D04-WT116A [Water] Sampled Oct-31-02 13:52 Central</b>				
GFAA 5100 Cd	Dec-03-02 12:00	30	Apr-29-03 13:52	
Volatiles Full List	Dec-03-02 12:00	30	Nov-07-02 13:52	
SVOA Standard List	Dec-03-02 12:00	30	Nov-07-02 13:52	
Cyanide, Total	Dec-03-02 12:00	30	Nov-14-02 13:52	pH10
Hg Total CVAA	Dec-03-02 12:00	30	Nov-28-02 13:52	pH2
Solids, Dry Weight	Dec-03-02 12:00	30	Nov-07-02 13:52	For SVOA Standard List in batch EK21101
B ICP (W)	Dec-03-02 12:00	30	Apr-29-03 13:52	pH2
THF & 1,4-Dioxane	Dec-03-02 12:00	30	Nov-14-02 13:52	
Metals SF ICP (W)	Dec-03-02 12:00	30	Apr-29-03 13:52	pH2
GFAA SIMAA Metals	Dec-03-02 12:00	30	Apr-29-03 13:52	pH2
<b>E2K0101-06 2003SY01S05-WT115A [Water] Sampled Oct-31-02 15:00 Central MS/MSD</b>				
GFAA 5100 Cd	Dec-03-02 12:00	30	Apr-29-03 15:00	
Volatiles Full List	Dec-03-02 12:00	30	Nov-07-02 15:00	
SVOA Standard List	Dec-03-02 12:00	30	Nov-07-02 15:00	
Cyanide, Total	Dec-03-02 12:00	30	Nov-14-02 15:00	pH10
Hg Total CVAA	Dec-03-02 12:00	30	Nov-28-02 15:00	pH2
Solids, Dry Weight	Dec-03-02 12:00	30	Nov-07-02 15:00	For SVOA Standard List in batch EK21101
B ICP (W)	Dec-03-02 12:00	30	Apr-29-03 15:00	pH2
THF & 1,4-Dioxane	Dec-03-02 12:00	30	Nov-14-02 15:00	
Metals SF ICP (W)	Dec-03-02 12:00	30	Apr-29-03 15:00	pH2
GFAA SIMAA Metals	Dec-03-02 12:00	30	Apr-29-03 15:00	pH2



**IIT RESEARCH INSTITUTE**

**IIT Research Institute ESAT Region 5**

**536 South Clark Street, Suite 734; Chicago, IL 60605**

**Telephone (312) 353-8302 Facsimile (312) 353-8307**

Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project:Himco Dump  
Project Number:2003SY01  
Project Manager:Howard Pham

Reported:  
Dec-05-02 14:50

**ANALYTICAL REPORT FOR SAMPLES**

Sample ID	Laboratory ID	Matrix	Date Sampled	Date Received
2003SY01R02-Pump Blank	E2K0101-02	Water	Oct-31-02 11:50	Nov-01-02 11:27
2003SY01S04-WT116A	E2K0101-04	Water	Oct-31-02 13:52	Nov-01-02 11:27
2003SY01D04-WT116A	E2K0101-05	Water	Oct-31-02 13:52	Nov-01-02 11:27
2003SY01S05-WT115A	E2K0101-06	Water	Oct-31-02 15:00	Nov-01-02 11:27



Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham

Reported:  
Dec-05-02 14:50

### 2003SY01R02-Pump Blank

E2K0101-02(Water)

Sampled: Oct-31-02 11:50

Received: Nov-01-02 11:27

#### Semivolatiles by GC/MS

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Phenol	U		0.74	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Bis(2-chloroethyl)ether	U		0.68	5.0	"	"	"	"	"
2-Chlorophenol	U		0.79	5.0	"	"	"	"	"
1,3-Dichlorobenzene	U		2.1	5.0	"	"	"	"	"
1,4-Dichlorobenzene	U		2.2	5.0	"	"	"	"	"
Benzyl alcohol	U		1.0	5.0	"	"	"	"	"
1,2-Dichlorobenzene	U		1.9	5.0	"	"	"	"	"
2-Methylphenol	U		0.64	5.0	"	"	"	"	"
Bis(2-chloroisopropyl)ether	U		1.8	5.0	"	"	"	"	"
4-Methylphenol	U		0.80	5.0	"	"	"	"	"
N-Nitrosodi-n-propylamine	U		1.7	5.0	"	"	"	"	"
Hexachloroethane	U		2.7	5.0	"	"	"	"	"
Nitrobenzene	U		1.4	5.0	"	"	"	"	"
Isophorone	U		1.4	5.0	"	"	"	"	"
2-Nitrophenol	U		1.5	5.0	"	"	"	"	"
2,4-Dimethylphenol	U		1.4	5.0	"	"	"	"	"
Benzoic acid	U	J	1.8	25	"	"	"	"	"
Bis(2-chloroethoxy)methane	U		1.0	5.0	"	"	"	"	"
2,4-Dichlorophenol	U		1.2	5.0	"	"	"	"	"
4-Trichlorobenzene	U		1.8	5.0	"	"	"	"	"
Naphthalene	U		1.1	5.0	"	"	"	"	"
4-Chloroaniline	U		1.6	5.0	"	"	"	"	"
Hexachlorobutadiene	U		2.5	5.0	"	"	"	"	"
2-Methylnaphthalene	U		1.1	5.0	"	"	"	"	"
4-Chloro-3-methylphenol	U		1.1	5.0	"	"	"	"	"
Hexachlorocyclopentadiene	U	J	2.8	25	"	"	"	"	"
2,4,6-Trichlorophenol	U		1.1	5.0	"	"	"	"	"
2,4,5-Trichlorophenol	U		1.6	5.0	"	"	"	"	"
2-Chloronaphthalene	U		1.4	5.0	"	"	"	"	"
2-Nitroaniline	U		2.3	5.0	"	"	"	"	"
Acenaphthylene	U		0.68	5.0	"	"	"	"	"
Dimethyl phthalate	U		1.2	5.0	"	"	"	"	"

**IIT RESEARCH INSTITUTE****IIT Research Institute ESAT Region 5****536 South Clark Street, Suite 734; Chicago, IL 60605****Telephone (312) 353-8302 Facsimile (312) 353-8307**Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham**Reported:**  
Dec-05-02 14:50**2003SY01R02-Pump Blank****E2K0101-02(Water)****Sampled: Oct-31-02 11:50****Received: Nov-01-02 11:27****Semivolatiles by GC/MS**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
2,6-Dinitrotoluene	U		1.6	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Acenaphthene	U		0.60	5.0	"	"	"	"	"
3-Nitroaniline	U		2.0	25	"	"	"	"	"
Dibenzofuran	U		1.6	5.0	"	"	"	"	"
2,4-Dinitrophenol	U	R	1.8	25	"	"	"	"	"
2,4-Dinitrotoluene	U		2.0	5.0	"	"	"	"	"
Fluorene	U		1.1	5.0	"	"	"	"	"
4-Nitrophenol	U		1.6	25	"	"	"	"	"
4-Chlorophenyl phenyl ether	U		1.4	5.0	"	"	"	"	"
Diethyl phthalate	U	J	1.7	5.0	"	"	"	"	"
4,6-Dinitro-2-methylphenol	U	J	1.5	25	"	"	"	"	"
N-Nitrosodiphenylamine	U		1.4	5.0	"	"	"	"	"
4-Nitroaniline	U		1.7	25	"	"	"	"	"
4-Bromophenyl phenyl ether	U		0.88	5.0	"	"	"	"	"
Hexachlorobenzene	U		1.4	5.0	"	"	"	"	"
Pentachlorophenol	U	J	1.6	25	"	"	"	"	"
Phenanthrene	U		0.56	5.0	"	"	"	"	"
Anthracene	U		0.80	5.0	"	"	"	"	"
Carbazole	U		1.6	5.0	"	"	"	"	"
Di-n-butyl phthalate	U		1.9	5.0	"	"	"	"	"
Fluoranthene	U		3.6	5.0	"	"	"	"	"
Pyrene	U		3.0	5.0	"	"	"	"	"
Butyl benzyl phthalate	U		2.2	5.0	"	"	"	"	"
3,3'-Dichlorobenzidine	U		1.1	25	"	"	"	"	"
Chrysene	U		0.65	5.0	"	"	"	"	"
Benzo (a) anthracene	U		0.92	5.0	"	"	"	"	"
Bis(2-ethylhexyl)phthalate	4.9	UJ	2.3	5.0	"	"	"	"	"
Di-n-octyl phthalate	U		2.7	5.0	"	"	"	"	"
Benzo (b) fluoranthene	U		1.1	5.0	"	"	"	"	"
Benzo (k) fluoranthene	U		1.1	5.0	"	"	"	"	"
Benzo (a) pyrene	U		0.73	5.0	"	"	"	"	"
Indeno (1,2,3-cd) pyrene	U		0.30	5.0	"	"	"	"	"



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77 West Jackson Boulevard  
Chicago IL, 60604

Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham

Reported:  
Dec-05-02 14:50

**2003SY01R02-Pump Blank**

**E2K0101-02(Water)**

**Sampled: Oct-31-02 11:50**

**Received: Nov-01-02 11:27**

**Semivolatiles by GC/MS**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dibenz (a,h) anthracene	U		0.63	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Benzo (g,h,i) perylene	U		0.93	5.0	"	"	"	"	"
Surrogate: 2-Fluorophenol	17.9			35.8 %		18-128	"	"	"
Surrogate: Phenol-d6	26.8			53.6 %		12-133	"	"	"
Surrogate: Nitrobenzene-d5	23.2			46.4 %		5-158	"	"	"
Surrogate: 2-Fluorobiphenyl	25.7			51.4 %		29-128	"	"	"
Surrogate: 2,4,6-Tribromophenol	39.0			78.0 %		32-148	"	"	"
Surrogate: Terphenyl-d14	42.3			84.6 %		62-140	"	"	"

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Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham

Reported:  
Dec-05-02 14:50

**2003SY01S04-WT116A****E2K0101-04(Water)****Sampled: Oct-31-02 13:52****Received: Nov-01-02 11:27****Semivolatiles by GC/MS**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Phenol	U		0.74	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Bis(2-chloroethyl)ether	U		0.68	5.0	"	"	"	"	"
2-Chlorophenol	U		0.79	5.0	"	"	"	"	"
1,3-Dichlorobenzene	U		2.1	5.0	"	"	"	"	"
1,4-Dichlorobenzene	U		2.2	5.0	"	"	"	"	"
Benzyl alcohol	U		1.0	5.0	"	"	"	"	"
1,2-Dichlorobenzene	U		1.9	5.0	"	"	"	"	"
2-Methylphenol	U		0.64	5.0	"	"	"	"	"
Bis(2-chloroisopropyl)ether	U		1.8	5.0	"	"	"	"	"
4-Methylphenol	U		0.80	5.0	"	"	"	"	"
N-Nitrosodi-n-propylamine	U		1.7	5.0	"	"	"	"	"
Hexachloroethane	U		2.7	5.0	"	"	"	"	"
Nitrobenzene	U		1.4	5.0	"	"	"	"	"
Isophorone	U		1.4	5.0	"	"	"	"	"
2-Nitrophenol	U		1.5	5.0	"	"	"	"	"
2,4-Dimethylphenol	U		1.4	5.0	"	"	"	"	"
Benzoic acid	U	J	1.8	25	"	"	"	"	"
Bis(2-chloroethoxy)methane	U		1.0	5.0	"	"	"	"	"
2,4-Dichlorophenol	U		1.2	5.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U		1.8	5.0	"	"	"	"	"
Naphthalene	U		1.1	5.0	"	"	"	"	"
4-Chloroaniline	U		1.6	5.0	"	"	"	"	"
Hexachlorobutadiene	U		2.5	5.0	"	"	"	"	"
2-Methylnaphthalene	U		1.1	5.0	"	"	"	"	"
4-Chloro-3-methylphenol	U		1.1	5.0	"	"	"	"	"
Hexachlorocyclopentadiene	U	J	2.8	25	"	"	"	"	"
2,4,6-Trichlorophenol	U		1.1	5.0	"	"	"	"	"
2,4,5-Trichlorophenol	U		1.6	5.0	"	"	"	"	"
2-Chloronaphthalene	U		1.4	5.0	"	"	"	"	"
2-Nitroaniline	U		2.3	5.0	"	"	"	"	"
Acenaphthylene	U		0.68	5.0	"	"	"	"	"
Dimethyl phthalate	U		1.2	5.0	"	"	"	"	"

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77 West Jackson Boulevard  
Chicago IL, 60604Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham**Reported:**  
Dec-05-02 14:50**2003SY01S04-WT116A****E2K0101-04(Water)****Sampled: Oct-31-02 13:52****Received: Nov-01-02 11:27****Semivolatiles by GC/MS**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
2,6-Dinitrotoluene	U		1.6	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Acenaphthene	U		0.60	5.0	"	"	"	"	"
3-Nitroaniline	U		2.0	25	"	"	"	"	"
2-Benzofuran	U		1.6	5.0	"	"	"	"	"
2,4-Dinitrophenol	U	R	1.8	25	"	"	"	"	"
2,4-Dinitrotoluene	U		2.0	5.0	"	"	"	"	"
Fluorene	U		1.1	5.0	"	"	"	"	"
4-Nitrophenol	U		1.6	25	"	"	"	"	"
4-Chlorophenyl phenyl ether	U		1.4	5.0	"	"	"	"	"
Diethyl phthalate	U	J	1.7	5.0	"	"	"	"	"
4,6-Dinitro-2-methylphenol	U	J	1.5	25	"	"	"	"	"
N-Nitrosodiphenylamine	U		1.4	5.0	"	"	"	"	"
4-Nitroaniline	U		1.7	25	"	"	"	"	"
4-Bromophenyl phenyl ether	U		0.88	5.0	"	"	"	"	"
Hexachlorobenzene	U		1.4	5.0	"	"	"	"	"
Pentachlorophenol	U	J	1.6	25	"	"	"	"	"
Phenanthrene	U		0.56	5.0	"	"	"	"	"
Anthracene	U		0.80	5.0	"	"	"	"	"
Carbazole	U		1.6	5.0	"	"	"	"	"
n-butyl phthalate	U		1.9	5.0	"	"	"	"	"
Fluoranthene	U		3.6	5.0	"	"	"	"	"
Pyrene	U		3.0	5.0	"	"	"	"	"
Butyl benzyl phthalate	U		2.2	5.0	"	"	"	"	"
3,3'-Dichlorobenzidine	U		1.1	25	"	"	"	"	"
Chrysene	U		0.65	5.0	"	"	"	"	"
Benzo (a) anthracene	U		0.92	5.0	"	"	"	"	"
Bis(2-ethylhexyl)phthalate	6.9	UJ	2.3	5.0	"	"	"	"	"
Di-n-octyl phthalate	U		2.7	5.0	"	"	"	"	"
Benzo (b) fluoranthene	U		1.1	5.0	"	"	"	"	"
Benzo (k) fluoranthene	U		1.1	5.0	"	"	"	"	"
Benzo (a) pyrene	U		0.73	5.0	"	"	"	"	"
Indeno (1,2,3-cd) pyrene	U		0.30	5.0	"	"	"	"	"



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Chicago IL, 60604

Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham

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Dec-05-02 14:50

2003SY01S04-WT116A

E2K0101-04(Water)

Sampled: Oct-31-02 13:52

Received: Nov-01-02 11:27

Semivolatiles by GC/MS

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dibenz (a,h) anthracene	U		0.63	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Benzo (g,h,i) perylene	U		0.93	5.0	"	"	"	"	"
Surrogate: 2-Fluorophenol	32.0			64.0 %		18-128	"	"	"
Surrogate: Phenol-d6	37.0			74.0 %		12-133	"	"	"
Surrogate: Nitrobenzene-d5	36.6			73.2 %		5-158	"	"	"
Surrogate: 2-Fluorobiphenyl	32.8			65.6 %		29-128	"	"	"
Surrogate: 2,4,6-Tribromophenol	52.4			105 %		32-148	"	"	"
Surrogate: Terphenyl-d14	39.0			78.0 %		62-140	"	"	"


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Project: Himco Dump  
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 Project Manager: Howard Pham

**Reported:**  
 Dec-05-02 14:50

**2003SY01D04-WT116A**

**E2K0101-05(Water)**

**Sampled: Oct-31-02 13:52**

**Received: Nov-01-02 11:27**

**Semivolatiles by GC/MS**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Phenol	U		0.74	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Bis(2-chloroethyl)ether	U		0.68	5.0	"	"	"	"	"
2-Chlorophenol	U		0.79	5.0	"	"	"	"	"
1,3-Dichlorobenzene	U		2.1	5.0	"	"	"	"	"
1,4-Dichlorobenzene	U		2.2	5.0	"	"	"	"	"
Benzyl alcohol	U		1.0	5.0	"	"	"	"	"
1,2-Dichlorobenzene	U		1.9	5.0	"	"	"	"	"
2-Methylphenol	U		0.64	5.0	"	"	"	"	"
Bis(2-chloroisopropyl)ether	U		1.8	5.0	"	"	"	"	"
4-Methylphenol	U		0.80	5.0	"	"	"	"	"
N-Nitrosodi-n-propylamine	U		1.7	5.0	"	"	"	"	"
Hexachloroethane	U		2.7	5.0	"	"	"	"	"
Nitrobenzene	U		1.4	5.0	"	"	"	"	"
Isophorone	U		1.4	5.0	"	"	"	"	"
2-Nitrophenol	U		1.5	5.0	"	"	"	"	"
2,4-Dimethylphenol	U		1.4	5.0	"	"	"	"	"
Benzoic acid	U	J	1.8	25	"	"	"	"	"
Bis(2-chloroethoxy)methane	U		1.0	5.0	"	"	"	"	"
2,4-Dichlorophenol	U		1.2	5.0	"	"	"	"	"
1,4-Trichlorobenzene	U		1.8	5.0	"	"	"	"	"
1-Naphthalene	U		1.1	5.0	"	"	"	"	"
4-Chloroaniline	U		1.6	5.0	"	"	"	"	"
Hexachlorobutadiene	U		2.5	5.0	"	"	"	"	"
2-Methylnaphthalene	U		1.1	5.0	"	"	"	"	"
4-Chloro-3-methylphenol	U		1.1	5.0	"	"	"	"	"
Hexachlorocyclopentadiene	U	J	2.8	25	"	"	"	"	"
2,4,6-Trichlorophenol	U		1.1	5.0	"	"	"	"	"
2,4,5-Trichlorophenol	U		1.6	5.0	"	"	"	"	"
2-Chloronaphthalene	U		1.4	5.0	"	"	"	"	"
2-Nitroaniline	U		2.3	5.0	"	"	"	"	"
Acenaphthylene	U		0.68	5.0	"	"	"	"	"
Dimethyl phthalate	U		1.2	5.0	"	"	"	"	"

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Project Manager: Howard Pham

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2003SY01D04-WT116A

E2K0101-05(Water)

Sampled: Oct-31-02 13:52

Received: Nov-01-02 11:27

#### Semivolatiles by GC/MS

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
2,6-Dinitrotoluene	U		1.6	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Acenaphthene	U		0.60	5.0	"	"	"	"	"
3-Nitroaniline	U		2.0	25	"	"	"	"	"
Dibenzofuran	U		1.6	5.0	"	"	"	"	"
2,4-Dinitrophenol	U	R	1.8	25	"	"	"	"	"
2,4-Dinitrotoluene	U		2.0	5.0	"	"	"	"	"
Fluorene	U		1.1	5.0	"	"	"	"	"
4-Nitrophenol	U		1.6	25	"	"	"	"	"
4-Chlorophenyl phenyl ether	U		1.4	5.0	"	"	"	"	"
Diethyl phthalate	U	J	1.7	5.0	"	"	"	"	"
4,6-Dinitro-2-methylphenol	U	J	1.5	25	"	"	"	"	"
N-Nitrosodiphenylamine	U		1.4	5.0	"	"	"	"	"
4-Nitroaniline	U		1.7	25	"	"	"	"	"
4-Bromophenyl phenyl ether	U		0.88	5.0	"	"	"	"	"
Hexachlorobenzene	U		1.4	5.0	"	"	"	"	"
Pentachlorophenol	U	J	1.6	25	"	"	"	"	"
Phenanthrene	U		0.56	5.0	"	"	"	"	"
Anthracene	U		0.80	5.0	"	"	"	"	"
Carbazole	U		1.6	5.0	"	"	"	"	"
Di-n-butyl phthalate	U		1.9	5.0	"	"	"	"	"
Fluoranthene	U		3.6	5.0	"	"	"	"	"
Pyrene	U		3.0	5.0	"	"	"	"	"
Butyl benzyl phthalate	U		2.2	5.0	"	"	"	"	"
3,3'-Dichlorobenzidine	U		1.1	25	"	"	"	"	"
Chrysene	U		0.65	5.0	"	"	"	"	"
Benzo (a) anthracene	U		0.92	5.0	"	"	"	"	"
Bis(2-ethylhexyl)phthalate	4.6	UJ	2.3	5.0	"	"	"	"	"
Di-n-octyl phthalate	U		2.7	5.0	"	"	"	"	"
Benzo (b) fluoranthene	U		1.1	5.0	"	"	"	"	"
Benzo (k) fluoranthene	U		1.1	5.0	"	"	"	"	"
Benzo (a) pyrene	U		0.73	5.0	"	"	"	"	"
Indeno (1,2,3-cd) pyrene	U		0.30	5.0	"	"	"	"	"



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77 West Jackson Boulevard  
Chicago IL, 60604Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham**Reported:**  
Dec-05-02 14:50**2003SY01D04-WT116A****E2K0101-05(Water)****Sampled: Oct-31-02 13:52****Received: Nov-01-02 11:27****Semivolatiles by GC/MS**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
<b>Dibenz (a,h) anthracene</b>	U		0.63	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
<b>Benzo (g,h,i) perylene</b>	U		0.93	5.0	"	"	"	"	"
<i>Surrogate: 2-Fluorophenol</i>	36.5			73.0 %		18-128	"	"	"
<i>Surrogate: Phenol-d6</i>	40.8			81.6 %		12-133	"	"	"
<i>Surrogate: Nitrobenzene-d5</i>	39.8			79.6 %		5-158	"	"	"
<i>Surrogate: 2-Fluorobiphenyl</i>	35.4			70.8 %		29-128	"	"	"
<i>Surrogate: 2,4,6-Tribromophenol</i>	53.0			106 %		32-148	"	"	"
<i>Surrogate: Terphenyl-d14</i>	42.4			84.8 %		62-140	"	"	"

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77 West Jackson Boulevard  
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Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham

Reported:  
Dec-05-02 14:50

2003SY01S05-WT115A

E2K0101-06(Water)

Sampled: Oct-31-02 15:00

Received: Nov-01-02 11:27

#### Semivolatiles by GC/MS

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Phenol	U		0.74	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Bis(2-chloroethyl)ether	U		0.68	5.0	"	"	"	"	"
2-Chlorophenol	U		0.79	5.0	"	"	"	"	"
1,3-Dichlorobenzene	U		2.1	5.0	"	"	"	"	"
1,4-Dichlorobenzene	U		2.2	5.0	"	"	"	"	"
Benzyl alcohol	U		1.0	5.0	"	"	"	"	"
1,2-Dichlorobenzene	U		1.9	5.0	"	"	"	"	"
2-Methylphenol	U		0.64	5.0	"	"	"	"	"
Bis(2-chloroisopropyl)ether	U		1.8	5.0	"	"	"	"	"
4-Methylphenol	U		0.80	5.0	"	"	"	"	"
N-Nitrosodi-n-propylamine	U		1.7	5.0	"	"	"	"	"
Hexachloroethane	U		2.7	5.0	"	"	"	"	"
Nitrobenzene	U		1.4	5.0	"	"	"	"	"
Isophorone	U		1.4	5.0	"	"	"	"	"
2-Nitrophenol	U		1.5	5.0	"	"	"	"	"
2,4-Dimethylphenol	U		1.4	5.0	"	"	"	"	"
Benzoic acid	U	J	1.8	25	"	"	"	"	"
Bis(2-chloroethoxy)methane	U		1.0	5.0	"	"	"	"	"
2,4-Dichlorophenol	U		1.2	5.0	"	"	"	"	"
1,2,4-Trichlorobenzene	U		1.8	5.0	"	"	"	"	"
Naphthalene	U		1.1	5.0	"	"	"	"	"
4-Chloroaniline	U		1.6	5.0	"	"	"	"	"
Hexachlorobutadiene	U		2.5	5.0	"	"	"	"	"
2-Methylnaphthalene	U		1.1	5.0	"	"	"	"	"
4-Chloro-3-methylphenol	U		1.1	5.0	"	"	"	"	"
Hexachlorocyclopentadiene	U	J	2.8	25	"	"	"	"	"
2,4,6-Trichlorophenol	U		1.1	5.0	"	"	"	"	"
2,4,5-Trichlorophenol	U		1.6	5.0	"	"	"	"	"
2-Chloronaphthalene	U		1.4	5.0	"	"	"	"	"
2-Nitroaniline	U		2.3	5.0	"	"	"	"	"
Acenaphthylene	U		0.68	5.0	"	"	"	"	"
Dimethyl phthalate	U		1.2	5.0	"	"	"	"	"

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Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham

**Reported:**  
Dec-05-02 14:50

**2003SY01S05-WT115A**

**E2K0101-06(Water)**

**Sampled: Oct-31-02 15:00**

**Received: Nov-01-02 11:27**

**Semivolatiles by GC/MS**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
2,6-Dinitrotoluene	U		1.6	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Acenaphthene	U		0.60	5.0	"	"	"	"	"
3-Nitroaniline	U		2.0	25	"	"	"	"	"
Dibenzofuran	U		1.6	5.0	"	"	"	"	"
m-Dinitrophenol	U	R	1.8	25	"	"	"	"	"
2,4-Dinitrotoluene	U		2.0	5.0	"	"	"	"	"
Fluorene	U		1.1	5.0	"	"	"	"	"
4-Nitrophenol	U		1.6	25	"	"	"	"	"
4-Chlorophenyl phenyl ether	U		1.4	5.0	"	"	"	"	"
Diethyl phthalate	U	J	1.7	5.0	"	"	"	"	"
4,6-Dinitro-2-methylphenol	U	J	1.5	25	"	"	"	"	"
N-Nitrosodiphenylamine	U		1.4	5.0	"	"	"	"	"
4-Nitroaniline	U		1.7	25	"	"	"	"	"
4-Bromophenyl phenyl ether	U		0.88	5.0	"	"	"	"	"
Hexachlorobenzene	U		1.4	5.0	"	"	"	"	"
Pentachlorophenol	U	J	1.6	25	"	"	"	"	"
Phenanthrene	U		0.56	5.0	"	"	"	"	"
Anthracene	U		0.80	5.0	"	"	"	"	"
Carbazole	U		1.6	5.0	"	"	"	"	"
n-butyl phthalate	U		1.9	5.0	"	"	"	"	"
Fluoranthene	U		3.6	5.0	"	"	"	"	"
Pyrene	U		3.0	5.0	"	"	"	"	"
Butyl benzyl phthalate	U		2.2	5.0	"	"	"	"	"
3,3'-Dichlorobenzidine	U	J	1.1	25	"	"	"	"	"
Chrysene	U		0.65	5.0	"	"	"	"	"
Benzo (a) anthracene	U		0.92	5.0	"	"	"	"	"
Bis(2-ethylhexyl)phthalate	6.0	UJ	2.3	5.0	"	"	"	"	"
Di-n-octyl phthalate	U		2.7	5.0	"	"	"	"	"
Benzo (b) fluoranthene	U		1.1	5.0	"	"	"	"	"
Benzo (k) fluoranthene	U		1.1	5.0	"	"	"	"	"
Benzo (a) pyrene	U		0.73	5.0	"	"	"	"	"
Indeno (1,2,3-cd) pyrene	U		0.30	5.0	"	"	"	"	"



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Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham

Reported:  
Dec-05-02 14:50

2003SY01S05-WT115A

E2K0101-06(Water)

Sampled: Oct-31-02 15:00

Received: Nov-01-02 11:27

Semivolatiles by GC/MS

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Dilution	Batch	Prepared	Analyzed
Dibenz (a,h) anthracene	U		0.63	5.0	ug/L	1	EK21101	Nov-04-02	Nov-20-02
Benzo (g,h,i) perylene	U		0.93	5.0	"	"	"	"	"
Surrogate: 2-Fluorophenol	27.8			55.6 %		18-128	"	"	"
Surrogate: Phenol-d6	35.6			71.2 %		12-133	"	"	"
Surrogate: Nitrobenzene-d5	26.7			53.4 %		5-158	"	"	"
Surrogate: 2-Fluorobiphenyl	30.3			60.6 %		29-128	"	"	"
Surrogate: 2,4,6-Tribromophenol	47.2			94.4 %		32-148	"	"	"
Surrogate: Terphenyl-d14	43.3			86.6 %		62-140	"	"	"

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Dec-05-02 14:50

**Semivolatiles by GC/MS - Quality Control**

Batch EK21101 - BP SVOA

Blank (EK21101-BLK1)

Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Phenol	U		0.74	5.0	ug/L					
Bis(2-chloroethyl)ether	U		0.68	5.0	"					
2-Chlorophenol	U		0.79	5.0	"					
1,3-Dichlorobenzene	U		2.1	5.0	"					
-Dichlorobenzene	U		2.2	5.0	"					
Benzyl alcohol	U		1.0	5.0	"					
1,2-Dichlorobenzene	U		1.9	5.0	"					
2-Methylphenol	U		0.64	5.0	"					
Bis(2-chloroisopropyl)ether	U		1.8	5.0	"					
4-Methylphenol	U		0.80	5.0	"					
N-Nitrosodi-n-propylamine	U		1.7	5.0	"					
Hexachloroethane	U		2.7	5.0	"					
Nitrobenzene	U		1.4	5.0	"					
Isophorone	U		1.4	5.0	"					
2-Nitrophenol	U		1.5	5.0	"					
2,4-Dimethylphenol	U		1.4	5.0	"					
Benzoic acid	U	J	1.8	25	"					
Bis(2-chloroethoxy)methane	U		1.0	5.0	"					
2,4-Dichlorophenol	U		1.2	5.0	"					
,4-Trichlorobenzene	U		1.8	5.0	"					
Naphthalene	U		1.1	5.0	"					
4-Chloroaniline	U		1.6	5.0	"					
Hexachlorobutadiene	U		2.5	5.0	"					
2-Methylnaphthalene	U		1.1	5.0	"					
4-Chloro-3-methylphenol	U		1.1	5.0	"					
Hexachlorocyclopentadiene	U	J	2.8	25	"					
2,4,6-Trichlorophenol	U		1.1	5.0	"					
2,4,5-Trichlorophenol	U		1.6	5.0	"					
2-Chloronaphthalene	U		1.4	5.0	"					
2-Nitroaniline	U		2.3	5.0	"					
Acenaphthylene	U		0.68	5.0	"					
Dimethyl phthalate	U		1.2	5.0	"					

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: Himco Dump  
 Project Number: 2003SY01  
 Project Manager: Howard Pham

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### Semivolatiles by GC/MS - Quality Control

#### Batch EK21101 - BP SVOA

Blank (EK21101-BLK1)

Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
2,6-Dinitrotoluene	U		1.6	5.0	ug/L					
Acenaphthene	U		0.60	5.0	"					
3-Nitroaniline	U		2.0	25	"					
Dibenzofuran	U		1.6	5.0	"					
2,4-Dinitrophenol	U	R	1.8	25	"					
2,4-Dinitrotoluene	U		2.0	5.0	"					
Fluorene	U		1.1	5.0	"					
4-Nitrophenol	U		1.6	25	"					
4-Chlorophenyl phenyl ether	U		1.4	5.0	"					
Diethyl phthalate	U		1.7	5.0	"					
4,6-Dinitro-2-methylphenol	U	J	1.5	25	"					
N-Nitrosodiphenylamine	U		1.4	5.0	"					
4-Nitroaniline	U		1.7	25	"					
4-Bromophenyl phenyl ether	U		0.88	5.0	"					
Hexachlorobenzene	U		1.4	5.0	"					
Pentachlorophenol	U	J	1.6	25	"					
Phenanthrene	U		0.56	5.0	"					
Anthracene	U		0.80	5.0	"					
Carbazole	U		1.6	5.0	"					
Di-n-butyl phthalate	U		1.9	5.0	"					
Fluoranthene	U		3.6	5.0	"					
Pyrene	U		3.0	5.0	"					
Butyl benzyl phthalate	U		2.2	5.0	"					
3,3'-Dichlorobenzidine	U		1.1	25	"					
Chrysene	U		0.65	5.0	"					
Benzo (a) anthracene	U		0.92	5.0	"					
Bis(2-ethylhexyl)phthalate	5.69	J	2.3	5.0	"					
Di-n-octyl phthalate	U		2.7	5.0	"					
Benzo (b) fluoranthene	U		1.1	5.0	"					
Benzo (k) fluoranthene	U		1.1	5.0	"					
Benzo (a) pyrene	U		0.73	5.0	"					
Indeno (1,2,3-cd) pyrene	U		0.30	5.0	"					



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**Telephone (312) 353-8302 Facsimile (312) 353-8307**

Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: Himco Dump  
 Project Number: 2003SY01  
 Project Manager: Howard Pham

**Reported:**  
 Dec-05-02 14:50

## Semivolatiles by GC/MS - Quality Control

### Batch EK21101 - BP SVOA

#### Blank (EK21101-BLK1)

Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Dibenz (a,h) anthracene	U		0.63	5.0	ug/L					
Benzo (g,h,i) perylene	U		0.93	5.0	"					
Surrogate: 2-Fluorophenol	21.5				"	50.0		43.0	18-128	
Surrogate: Phenol-d6	29.8				"	50.0		59.6	12-133	
Surrogate: Nitrobenzene-d5	30.6				"	50.0		61.2	5-158	
Surrogate: 2-Fluorobiphenyl	28.0				"	50.0		56.0	29-128	
Surrogate: 2,4,6-Tribromophenol	37.1				"	50.0		74.2	32-148	
Surrogate: Terphenyl-d14	40.8				"	50.0		81.6	62-140	

#### LCS (EK21101-BS1)

Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Phenol	34.5		0.74	5.0	ug/L	50.0		69.0	7-128	
Bis(2-chloroethyl)ether	31.0		0.68	5.0	"	50.0		62.0	33-114	
2-Chlorophenol	35.2		0.79	5.0	"	50.0		70.4	31-121	
1,3-Dichlorobenzene	22.7		2.1	5.0	"	50.0		45.4	21-101	
1,4-Dichlorobenzene	24.6		2.2	5.0	"	50.0		49.2	22-102	
Benzyl alcohol	37.4		1.0	5.0	"	50.0		74.8	34-132	
1,2-Dichlorobenzene	26.6		1.9	5.0	"	50.0		53.2	27-102	
2-Methylphenol	33.6		0.64	5.0	"	50.0		67.2	36-115	
Bis(2-chloroisopropyl)ether	34.3		1.8	5.0	"	50.0		68.6	34-120	
4-Methylphenol	33.7		0.80	5.0	"	50.0		67.4	35-117	
N-Nitrosodi-n-propylamine	35.2		1.7	5.0	"	50.0		70.4	49-111	
Hexachloroethane	14.1		2.7	5.0	"	50.0		28.2	16-103	
Nitrobenzene	33.8		1.4	5.0	"	50.0		67.6	41-115	
Isophorone	36.1		1.4	5.0	"	50.0		72.2	51-115	
2-Nitrophenol	34.4		1.5	5.0	"	50.0		68.8	41-128	
2,4-Dimethylphenol	26.8		1.4	5.0	"	50.0		53.6	13-123	
Benzoic acid	U	J	1.8	25	"	50.0			5-176	
Bis(2-chloroethoxy)methane	34.1		1.0	5.0	"	50.0		68.2	48-118	
2,4-Dichlorophenol	36.8		1.2	5.0	"	50.0		73.6	40-137	

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### Semivolatiles by GC/MS - Quality Control

#### Batch EK21101 - BP SVOA

LCS (EK21101-BS1)

Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
1,2,4-Trichlorobenzene	26.9		1.8	5.0	ug/L	50.0		53.8	36-108	
Naphthalene	32.5		1.1	5.0	"	50.0		65.0	36-111	
4-Chloroaniline	38.0		1.6	5.0	"	50.0		76.0	50-133	
Hexachlorobutadiene	19.5		2.5	5.0	"	50.0		39.0	26-108	
2-Methylnaphthalene	31.2		1.1	5.0	"	50.0		62.4	47-111	
4-Chloro-3-methylphenol	38.6		1.1	5.0	"	50.0		77.2	52-131	
Hexachlorocyclopentadiene	4.25	J	2.8	25	"	50.0		8.50	5-179	
2,4,6-Trichlorophenol	41.1		1.1	5.0	"	50.0		82.2	38-139	
2,4,5-Trichlorophenol	39.7		1.6	5.0	"	50.0		79.4	34-143	
2-Chloronaphthalene	34.5		1.4	5.0	"	50.0		69.0	50-114	
2-Nitroaniline	44.2		2.3	5.0	"	50.0		88.4	55-131	
Acenaphthylene	36.4		0.68	5.0	"	50.0		72.8	44-121	
Dimethyl phthalate	3.57		1.2	5.0	"	50.0		7.14	4-102	
2,6-Dinitrotoluene	38.1		1.6	5.0	"	50.0		76.2	58-128	
Acenaphthene	36.2		0.60	5.0	"	50.0		72.4	52-114	
3-Nitroaniline	42.9		2.0	25	"	50.0		85.8	49-137	
Dibenzofuran	35.2		1.6	5.0	"	50.0		70.4	52-119	
2,4-Dinitrophenol	U	J	1.8	25	"	50.0			16-189	
2,4-Dinitrotoluene	40.1		2.0	5.0	"	50.0		80.2	57-132	
Fluorene	38.9		1.1	5.0	"	50.0		77.8	53-120	
4-Nitrophenol	70.6		1.6	25	"	50.0		141	5-235	
4-Chlorophenyl phenyl ether	37.8		1.4	5.0	"	50.0		75.6	54-124	
Diethyl phthalate	10.3		1.7	5.0	"	50.0		20.6	34-129	
4,6-Dinitro-2-methylphenol	17.6	J	1.5	25	"	50.0		35.2	11-235	
N-Nitrosodiphenylamine	34.3		1.4	5.0	"	50.0		68.6	42-154	
4-Nitroaniline	52.0		1.7	25	"	50.0		104	4-158	
4-Bromophenyl phenyl ether	36.3		0.88	5.0	"	50.0		72.6	49-148	
Hexachlorobenzene	38.2		1.4	5.0	"	50.0		76.4	52-143	
Pentachlorophenol	U	J	1.6	25	"	50.0			11-247	
Phenanthrene	38.1		0.56	5.0	"	50.0		76.2	44-144	
Anthracene	38.4		0.80	5.0	"	50.0		76.8	44-144	
Carbazole	43.0		1.6	5.0	"	50.0		86.0	49-139	





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Superfund, US EPA Region 5  
 77 West Jackson Boulevard  
 Chicago IL, 60604

Project: Himco Dump  
 Project Number: 2003SY01  
 Project Manager: Howard Pham

Reported:  
 Dec-05-02 14:50

**Semivolatiles by GC/MS - Quality Control**

**Batch EK21101 - BP SVOA**

**LCS (EK21101-BS1)**

**Prepared: Nov-04-02 Analyzed: Nov-20-02**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Di-n-butyl phthalate	37.0		1.9	5.0	ug/L	50.0		74.0	31-159		
Fluoranthene	42.5		3.6	5.0	"	50.0		85.0	49-138		
Pyrene	44.4		3.0	5.0	"	50.0		88.8	40-158		
Butyl benzyl phthalate	47.7		2.2	5.0	"	50.0		95.4	48-159		
1,3-Dichlorobenzidine	51.6		1.1	25	"	50.0		103	41-183		
Chrysene	38.6		0.65	5.0	"	50.0		77.2	50-152		
Benzo (a) anthracene	42.7		0.92	5.0	"	50.0		85.4	54-153		
Bis(2-ethylhexyl)phthalate	45.4	J	2.3	5.0	"	50.0		90.8	47-153		
Di-n-octyl phthalate	44.3		2.7	5.0	"	50.0		88.6	30-156		
Benzo (b) fluoranthene	41.6		1.1	5.0	"	50.0		83.2	55-153		
Benzo (k) fluoranthene	40.4		1.1	5.0	"	50.0		80.8	43-149		
Benzo (a) pyrene	43.2		0.73	5.0	"	50.0		86.4	47-156		
Indeno (1,2,3-cd) pyrene	36.9		0.30	5.0	"	50.0		73.8	43-153		
Dibenz (a,h) anthracene	36.4		0.63	5.0	"	50.0		72.8	48-146		
Benzo (g,h,i) perylene	33.8		0.93	5.0	"	50.0		67.6	37-155		
Surrogate: 2-Fluorophenol	35.4				"	50.0		70.8	18-128		
Surrogate: Phenol-d6	36.1				"	50.0		72.2	12-133		
Surrogate: Nitrobenzene-d5	37.1				"	50.0		74.2	5-158		
Surrogate: 2-Fluorobiphenyl	35.9				"	50.0		71.8	29-128		
Surrogate: 2,4,6-Tribromophenol	48.6				"	50.0		97.2	32-148		
Surrogate: Terphenyl-d14	48.6				"	50.0		97.2	62-140		

**LCS Dup (EK21101-BSD1)**

**Prepared: Nov-04-02 Analyzed: Nov-20-02**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Phenol	33.8		0.74	5.0	ug/L	50.0		67.6	7-128	2.05	84
Bis(2-chloroethyl)ether	31.2		0.68	5.0	"	50.0		62.4	33-114	0.643	63
2-Chlorophenol	34.9		0.79	5.0	"	50.0		69.8	31-121	0.856	62
1,3-Dichlorobenzene	24.4		2.1	5.0	"	50.0		48.8	21-101	7.22	75
1,4-Dichlorobenzene	25.7		2.2	5.0	"	50.0		51.4	22-102	4.37	79
Benzyl alcohol	36.0		1.0	5.0	"	50.0		72.0	34-132	3.81	40

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Project: Himco Dump  
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Project Manager: Howard Pham

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**Semivolatiles by GC/MS - Quality Control****Batch EK21101 - BP SVOA****LCS Dup (EK21101-BSD1)****Prepared: Nov-04-02 Analyzed: Nov-20-02**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
1,2-Dichlorobenzene	27.6		1.9	5.0	ug/L	50.0	55.2	27-102	3.69	73
2-Methylphenol	32.9		0.64	5.0	"	50.0	65.8	36-115	2.11	53
Bis(2-chloroisopropyl)ether	33.1		1.8	5.0	"	50.0	66.2	34-120	3.56	63
4-Methylphenol	32.4		0.80	5.0	"	50.0	64.8	35-117	3.93	51
N-Nitrosodi-n-propylamine	33.8		1.7	5.0	"	50.0	67.6	49-111	4.06	37
Hexachloroethane	16.4		2.7	5.0	"	50.0	32.8	16-103	15.1	77
Nitrobenzene	33.0		1.4	5.0	"	50.0	66.0	41-115	2.40	55
Isophorone	34.5		1.4	5.0	"	50.0	69.0	51-115	4.53	30
2-Nitrophenol	34.2		1.5	5.0	"	50.0	68.4	41-128	0.583	50
2,4-Dimethylphenol	27.8		1.4	5.0	"	50.0	55.6	13-123	3.66	73
Benzoic acid	U	J	1.8	25	"	50.0		5-176		155
Bis(2-chloroethoxy)methane	33.0		1.0	5.0	"	50.0	66.0	48-118	3.28	38
2,4-Dichlorophenol	36.2		1.2	5.0	"	50.0	72.4	40-137	1.64	53
1,2,4-Trichlorobenzene	26.8		1.8	5.0	"	50.0	53.6	36-108	0.372	58
Naphthalene	31.8		1.1	5.0	"	50.0	63.6	36-111	2.18	48
4-Chloroaniline	36.4		1.6	5.0	"	50.0	72.8	50-133	4.30	72
Hexachlorobutadiene	19.1		2.5	5.0	"	50.0	38.2	26-108	2.07	65
2-Methylnaphthalene	30.0		1.1	5.0	"	50.0	60.0	47-111	3.92	37
4-Chloro-3-methylphenol	37.0		1.1	5.0	"	50.0	74.0	52-131	4.23	37
Hexachlorocyclopentadiene	5.11	J	2.8	25	"	50.0	10.2	5-179	18.4	103
2,4,6-Trichlorophenol	39.9		1.1	5.0	"	50.0	79.8	38-139	2.96	48
2,4,5-Trichlorophenol	38.8		1.6	5.0	"	50.0	77.6	34-143	2.29	46
2-Chloronaphthalene	33.0		1.4	5.0	"	50.0	66.0	50-114	4.44	35
2-Nitroaniline	41.8		2.3	5.0	"	50.0	83.6	55-131	5.58	28
Acenaphthylene	34.5		0.68	5.0	"	50.0	69.0	44-121	5.36	29
Dimethyl phthalate	10.8		1.2	5.0	"	50.0	21.6	4-102	101	60
2,6-Dinitrotoluene	36.0		1.6	5.0	"	50.0	72.0	58-128	5.67	27
Acenaphthene	34.1		0.60	5.0	"	50.0	68.2	52-114	5.97	31
3-Nitroaniline	41.2		2.0	25	"	50.0	82.4	49-137	4.04	28
Dibenzofuran	33.7		1.6	5.0	"	50.0	67.4	52-119	4.35	27
2,4-Dinitrophenol	U	J	1.8	25	"	50.0		16-189		109
2,4-Dinitrotoluene	37.2		2.0	5.0	"	50.0	74.4	57-132	7.50	29

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Reported:  
Dec-05-02 14:50

### Semivolatiles by GC/MS - Quality Control

Batch EK21101 - BP SVOA

LCS Dup (EK21101-BSD1)

Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC Limits	RPD	RPD Limit
Fluorene	36.4		1.1	5.0	ug/L	50.0	72.8	53-120	6.64	27
4-Nitrophenol	68.2		1.6	25	"	50.0	136	5-235	3.46	64
4-Chlorophenyl phenyl ether	35.2		1.4	5.0	"	50.0	70.4	54-124	7.12	30
Diethyl phthalate	16.5		1.7	5.0	"	50.0	33.0	34-129	46.3	37
-Dinitro-2-methylphenol	17.2	J	1.5	25	"	50.0	34.4	11-235	2.30	59
N-Nitrosodiphenylamine	33.6		1.4	5.0	"	50.0	67.2	42-154	2.06	54
4-Nitroaniline	48.5		1.7	25	"	50.0	97.0	4-158	6.97	44
4-Bromophenyl phenyl ether	35.5		0.88	5.0	"	50.0	71.0	49-148	2.23	30
Hexachlorobenzene	36.3		1.4	5.0	"	50.0	72.6	52-143	5.10	29
Pentachlorophenol	4.96	J	1.6	25	"	50.0	9.92	11-247		87
Phenanthrene	36.0		0.56	5.0	"	50.0	72.0	44-144	5.67	25
Anthracene	36.3		0.80	5.0	"	50.0	72.6	44-144	5.62	29
Carbazole	40.5		1.6	5.0	"	50.0	81.0	49-139	5.99	47
Di-n-butyl phthalate	34.6		1.9	5.0	"	50.0	69.2	31-159	6.70	40
Fluoranthene	39.8		3.6	5.0	"	50.0	79.6	49-138	6.56	30
Pyrene	40.9		3.0	5.0	"	50.0	81.8	40-158	8.21	28
Butyl benzyl phthalate	43.0		2.2	5.0	"	50.0	86.0	48-159	10.4	40
3,3'-Dichlorobenzidine	50.0		1.1	25	"	50.0	100	41-183	3.15	44
Chrysene	35.4		0.65	5.0	"	50.0	70.8	50-152	8.65	36
Benzo (a) anthracene	39.8		0.92	5.0	"	50.0	79.6	54-153	7.03	36
Diethylhexyl phthalate	41.3	J	2.3	5.0	"	50.0	82.6	47-153	9.46	22
Di-n-octyl phthalate	38.4		2.7	5.0	"	50.0	76.8	30-156	14.3	87
Benzo (b) fluoranthene	85.8		1.1	5.0	"	50.0	172	55-153	69.4	83
Benzo (k) fluoranthene	73.9		1.1	5.0	"	50.0	148	43-149	58.6	83
Benzo (a) pyrene	39.8		0.73	5.0	"	50.0	79.6	47-156	8.19	72
Indeno (1,2,3-cd) pyrene	33.1		0.30	5.0	"	50.0	66.2	43-153	10.9	66
Dibenz (a,h) anthracene	33.2		0.63	5.0	"	50.0	66.4	48-146	9.20	74
Benzo (g,h,i) perylene	29.7		0.93	5.0	"	50.0	59.4	37-155	12.9	58
Surrogate: 2-Fluorophenol	35.0				"	50.0	70.0	18-128		
Surrogate: Phenol-d6	34.7				"	50.0	69.4	12-133		
Surrogate: Nitrobenzene-d5	34.9				"	50.0	69.8	5-158		
Surrogate: 2-Fluorobiphenyl	33.0				"	50.0	66.0	29-128		



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Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604

Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham

Reported:  
Dec-05-02 14:50

**Semivolatiles by GC/MS - Quality Control**

**Batch EK21101 - BP SVOA**

**LCS Dup (EK21101-BSD1)**

**Prepared: Nov-04-02 Analyzed: Nov-20-02**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Surrogate: 2,4,6-Tribromophenol	47.0				ug/L	50.0		94.0	32-148		
Surrogate: Terphenyl-d14	43.6				"	50.0		87.2	62-140		

**Matrix Spike (EK21101-MS1)**

**Source: E2K0101-06**

**Prepared: Nov-04-02 Analyzed: Nov-20-02**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Phenol	31.6		0.74	5.0	ug/L	50.0	U	63.2	5-131		
Bis(2-chloroethyl)ether	19.3		0.68	5.0	"	50.0	U	38.6	5-156		
2-Chlorophenol	31.1		0.79	5.0	"	50.0	U	62.2	5-138		
1,3-Dichlorobenzene	23.1		2.1	5.0	"	50.0	U	46.2	39-96		
1,4-Dichlorobenzene	23.9		2.2	5.0	"	50.0	U	47.8	36-98		
Benzyl alcohol	34.2		1.0	5.0	"	50.0	U	68.4	5-188		
1,2-Dichlorobenzene	25.2		1.9	5.0	"	50.0	U	50.4	40-99		
2-Methylphenol	31.5		0.64	5.0	"	50.0	U	63.0	21-117		
Bis(2-chloroisopropyl)ether	28.5		1.8	5.0	"	50.0	U	57.0	5-168		
4-Methylphenol	31.7		0.80	5.0	"	50.0	U	63.4	26-114		
N-Nitrosodi-n-propylamine	30.9		1.7	5.0	"	50.0	U	61.8	5-182		
Hexachloroethane	22.6		2.7	5.0	"	50.0	U	45.2	5-132		
Nitrobenzene	29.5		1.4	5.0	"	50.0	U	59.0	5-163		
Isophorone	32.4		1.4	5.0	"	50.0	U	64.8	5-171		
2-Nitrophenol	31.7		1.5	5.0	"	50.0	U	63.4	5-157		
2,4-Dimethylphenol	26.0		1.4	5.0	"	50.0	U	52.0	5-116		
Benzoic acid	U	J	1.8	25	"	50.0	U		5-150		
Bis(2-chloroethoxy)methane	30.8		1.0	5.0	"	50.0	U	61.6	5-174		
2,4-Dichlorophenol	34.4		1.2	5.0	"	50.0	U	68.8	24-127		
1,2,4-Trichlorobenzene	29.2		1.8	5.0	"	50.0	U	58.4	41-106		
Naphthalene	29.9		1.1	5.0	"	50.0	U	59.8	5-168		
4-Chloroaniline	28.9		1.6	5.0	"	50.0	U	57.8	5-183		
Hexachlorobutadiene	28.9		2.5	5.0	"	50.0	U	57.8	38-99		
2-Methylnaphthalene	30.7		1.1	5.0	"	50.0	U	61.4	48-109		
4-Chloro-3-methylphenol	38.4		1.1	5.0	"	50.0	U	76.8	27-140		

Xuyen Nguyen, Chemist

Report Name: E2K0101

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**IIT RESEARCH INSTITUTE****IIT Research Institute ESAT Region 5****536 South Clark Street, Suite 734; Chicago, IL 60605****Telephone (312) 353-8302 Facsimile (312) 353-8307**Superfund, US EPA Region 5  
77 West Jackson Boulevard  
Chicago IL, 60604Project: Himco Dump  
Project Number: 2003SY01  
Project Manager: Howard Pham**Reported:**  
Dec-05-02 14:50**Semivolatiles by GC/MS - Quality Control****Batch EK21101 - BP SVOA****Matrix Spike (EK21101-MS1)****Source: E2K0101-06****Prepared: Nov-04-02 Analyzed: Nov-20-02**

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Hexachlorocyclopentadiene	7.76	J	2.8	25	ug/L	50.0	U	15.5	5-95		
2,4,6-Trichlorophenol	41.0		1.1	5.0	"	50.0	U	82.0	26-132		
2,4,5-Trichlorophenol	39.5		1.6	5.0	"	50.0	U	79.0	47-123		
2-Chloronaphthalene	34.2		1.4	5.0	"	50.0	U	68.4	56-106		
Nitroaniline	41.5		2.3	5.0	"	50.0	U	83.0	5-176		
Acenaphthylene	34.8		0.68	5.0	"	50.0	U	69.6	5-174		
Dimethyl phthalate	35.7		1.2	5.0	"	50.0	U	71.4	39-102		
2,6-Dinitrotoluene	38.4		1.6	5.0	"	50.0	U	76.8	67-110		
Acenaphthene	35.2		0.60	5.0	"	50.0	U	70.4	18-129		
3-Nitroaniline	37.6		2.0	25	"	50.0	U	75.2	5-203		
Dibenzofuran	33.9		1.6	5.0	"	50.0	U	67.8	5-112		
2,4-Dinitrophenol	U	J	1.8	25	"	50.0	U		5-191		
2,4-Dinitrotoluene	40.2		2.0	5.0	"	50.0	U	80.4	28-146		
Fluorene	37.4		1.1	5.0	"	50.0	U	74.8	55-113		
4-Nitrophenol	82.6		1.6	25	"	50.0	U	165	34-128		
4-Chlorophenyl phenyl ether	36.9		1.4	5.0	"	50.0	U	73.8	59-113		
Diethyl phthalate	37.3		1.7	5.0	"	50.0	U	74.6	55-100		
4,6-Dinitro-2-methylphenol	30.5	J	1.5	25	"	50.0	U	61.0	72-138		
N-Nitrosodiphenylamine	35.8		1.4	5.0	"	50.0	U	71.6	74-102		
Nitroaniline	44.3		1.7	25	"	50.0	U	88.6	5-177		
Bromophenyl phenyl ether	36.4		0.88	5.0	"	50.0	U	72.8	71-107		
Hexachlorobenzene	37.9		1.4	5.0	"	50.0	U	75.8	76-100		
Pentachlorophenol	94.4	J	1.6	25	"	50.0	U	189	5-394		
Phenanthrene	37.8		0.56	5.0	"	50.0	U	75.6	65-112		
Anthracene	37.8		0.80	5.0	"	50.0	U	75.6	61-114		
Carbazole	42.2		1.6	5.0	"	50.0	U	84.4	50-146		
Di-n-butyl phthalate	37.5		1.9	5.0	"	50.0	U	75.0	51-126		
Fluoranthene	41.2		3.6	5.0	"	50.0	U	82.4	41-152		
Pyrene	42.1		3.0	5.0	"	50.0	U	84.2	57-119		
Butyl benzyl phthalate	45.8		2.2	5.0	"	50.0	U	91.6	20-152		
3,3'-Dichlorobenzidine	13.3		1.1	25	"	50.0	U	26.6	5-113		
Chrysene	37.0		0.65	5.0	"	50.0	U	74.0	62-118		

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## Semivolatiles by GC/MS - Quality Control

### Batch EK21101 - BP SVOA

Matrix Spike (EK21101-MS1) Source: E2K0101-06 Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Benzo (a) anthracene	42.0		0.92	5.0	ug/L	50.0	U	84.0	63-120		
Bis(2-ethylhexyl)phthalate	65.5	J	2.3	5.0	"	50.0	6.0	119	39-138		
Di-n-octyl phthalate	38.6		2.7	5.0	"	50.0	U	77.2	28-158		
Benzo (b) fluoranthene	45.1		1.1	5.0	"	50.0	U	90.2	49-155		
Benzo (k) fluoranthene	39.8		1.1	5.0	"	50.0	U	79.6	43-149		
Benzo (a) pyrene	41.5		0.73	5.0	"	50.0	U	83.0	31-160		
Indeno (1,2,3-cd) pyrene	38.9		0.30	5.0	"	50.0	U	77.8	52-171		
Dibenz (a,h) anthracene	38.9		0.63	5.0	"	50.0	U	77.8	12-226		
Benzo (g,h,i) perylene	37.0		0.93	5.0	"	50.0	U	74.0	25-204		
Surrogate: 2-Fluorophenol	30.4				"	50.0		60.8	18-128		
Surrogate: Phenol-d6	34.1				"	50.0		68.2	12-133		
Surrogate: Nitrobenzene-d5	32.1				"	50.0		64.2	5-158		
Surrogate: 2-Fluorobiphenyl	34.5				"	50.0		69.0	29-128		
Surrogate: 2,4,6-Tribromophenol	49.1				"	50.0		98.2	32-148		
Surrogate: Terphenyl-d14	44.5				"	50.0		89.0	62-140		

Matrix Spike Dup (EK21101-MSD1) Source: E2K0101-06 Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Phenol	27.2		0.74	5.0	ug/L	50.0	U	54.4	5-131	15.0	42
Bis(2-chloroethyl)ether	4.94		0.68	5.0	"	50.0	U	9.88	5-156	118	45
2-Chlorophenol	21.9		0.79	5.0	"	50.0	U	43.8	5-138	34.7	86
1,3-Dichlorobenzene	5.51		2.1	5.0	"	50.0	U	11.0	39-96	123	82
1,4-Dichlorobenzene	6.09		2.2	5.0	"	50.0	U	12.2	36-98	119	75
Benzyl alcohol	26.7		1.0	5.0	"	50.0	U	53.4	5-188	24.6	28
1,2-Dichlorobenzene	7.77		1.9	5.0	"	50.0	U	15.5	40-99	106	66
2-Methylphenol	26.4		0.64	5.0	"	50.0	U	52.8	21-117	17.6	62
Bis(2-chloroisopropyl)ether	11.6		1.8	5.0	"	50.0	U	23.2	5-168	84.3	39
4-Methylphenol	27.9		0.80	5.0	"	50.0	U	55.8	26-114	12.8	44
N-Nitrosodi-n-propylamine	16.5		1.7	5.0	"	50.0	U	33.0	5-182	60.8	48
Hexachloroethane	7.83		2.7	5.0	"	50.0	U	15.7	5-132	97.1	200

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Project Manager: Howard Pham

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## Semivolatiles by GC/MS - Quality Control

### Batch EK21101 - BP SVOA

Matrix Spike Dup (EK21101-MSD1) Source: E2K0101-06 Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
Nitrobenzene	15.6		1.4	5.0	ug/L	50.0	U	31.2	5-163	61.6	110
Isophorone	20.7		1.4	5.0	"	50.0	U	41.4	5-171	44.1	36
2-Nitrophenol	19.1		1.5	5.0	"	50.0	U	38.2	5-157	49.6	75
2,4-Dimethylphenol	13.6		1.4	5.0	"	50.0	U	27.2	5-116	62.6	69
Azoic acid	U	J	1.8	25	"	50.0	U		5-150		42
Bis(2-chloroethoxy)methane	21.8		1.0	5.0	"	50.0	U	43.6	5-174	34.2	296
2,4-Dichlorophenol	34.9		1.2	5.0	"	50.0	U	69.8	24-127	1.44	23
1,2,4-Trichlorobenzene	16.3		1.8	5.0	"	50.0	U	32.6	41-106	56.7	115
Naphthalene	18.8		1.1	5.0	"	50.0	U	37.6	5-168	45.6	37
4-Chloroaniline	12.7		1.6	5.0	"	50.0	U	25.4	5-183	77.9	30
Hexachlorobutadiene	15.1		2.5	5.0	"	50.0	U	30.2	38-99	62.7	145
2-Methylnaphthalene	24.6		1.1	5.0	"	50.0	U	49.2	48-109	22.1	34
4-Chloro-3-methylphenol	39.1		1.1	5.0	"	50.0	U	78.2	27-140	1.81	28
Hexachlorocyclopentadiene	5.40	J	2.8	25	"	50.0	U	10.8	5-95	35.9	168
2,4,6-Trichlorophenol	40.8		1.1	5.0	"	50.0	U	81.6	26-132	0.489	25
2,4,5-Trichlorophenol	41.8		1.6	5.0	"	50.0	U	83.6	47-123	5.66	24
2-Chloronaphthalene	32.2		1.4	5.0	"	50.0	U	64.4	56-106	6.02	26
2-Nitroaniline	43.6		2.3	5.0	"	50.0	U	87.2	5-176	4.94	37
Acenaphthylene	33.9		0.68	5.0	"	50.0	U	67.8	5-174	2.62	23
Dimethyl phthalate	37.4		1.2	5.0	"	50.0	U	74.8	39-102	4.65	44
2,6-Dinitrotoluene	39.4		1.6	5.0	"	50.0	U	78.8	67-110	2.57	20
Acenaphthene	35.1		0.60	5.0	"	50.0	U	70.2	18-129	0.285	54
3-Nitroaniline	35.6		2.0	25	"	50.0	U	71.2	5-203	5.46	28
Dibenzofuran	35.9		1.6	5.0	"	50.0	U	71.8	5-112	5.73	16
2,4-Dinitrophenol	19.2	J	1.8	25	"	50.0	U	38.4	5-191		96
2,4-Dinitrotoluene	41.3		2.0	5.0	"	50.0	U	82.6	28-146	2.70	19
Fluorene	39.4		1.1	5.0	"	50.0	U	78.8	55-113	5.21	21
4-Nitrophenol	113		1.6	25	"	50.0	U	226	34-128	31.1	200
4-Chlorophenyl phenyl ether	38.6		1.4	5.0	"	50.0	U	77.2	59-113	4.50	17
Diethyl phthalate	39.4		1.7	5.0	"	50.0	U	78.8	55-100	5.48	22
4,6-Dinitro-2-methylphenol	24.3	J	1.5	25	"	50.0	U	48.6	72-138	22.6	68
N-Nitrosodiphenylamine	36.4		1.4	5.0	"	50.0	U	72.8	74-102	1.66	31

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 Project Manager: Howard Pham

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### Semivolatiles by GC/MS - Quality Control

#### Batch EK21101 - BP SVOA

Matrix Spike Dup (EK21101-MSD1) Source: E2K0101-06 Prepared: Nov-04-02 Analyzed: Nov-20-02

Analyte	Result	Flags / Qualifiers	MDL	Reporting Limit	Units	Spike Level	Source Result	%REC	%REC Limits	RPD	RPD Limit
4-Nitroaniline	47.3		1.7	25	ug/L	50.0	U	94.6	5-177	6.55	73
4-Bromophenyl phenyl ether	37.9		0.88	5.0	"	50.0	U	75.8	71-107	4.04	12
Hexachlorobenzene	39.1		1.4	5.0	"	50.0	U	78.2	76-100	3.12	17
Pentachlorophenol	80.6	J	1.6	25	"	50.0	U	161	5-394	15.8	57
Phenanthrene	39.6		0.56	5.0	"	50.0	U	79.2	65-112	4.65	16
Anthracene	33.0		0.80	5.0	"	50.0	U	66.0	61-114	13.6	33
Carbazole	45.6		1.6	5.0	"	50.0	U	91.2	50-146	7.74	30
Di-n-butyl phthalate	39.6		1.9	5.0	"	50.0	U	79.2	51-126	5.45	40
Fluoranthene	44.2		3.6	5.0	"	50.0	U	88.4	41-152	7.03	22
Pyrene	44.5		3.0	5.0	"	50.0	U	89.0	57-119	5.54	55
Butyl benzyl phthalate	49.5		2.2	5.0	"	50.0	U	99.0	20-152	7.76	70
3,3'-Dichlorobenzidine	8.81		1.1	25	"	50.0	U	17.6	5-113	40.6	27
Chrysene	38.8		0.65	5.0	"	50.0	U	77.6	62-118	4.75	17
Benzo (a) anthracene	42.7		0.92	5.0	"	50.0	U	85.4	63-120	1.65	23
Bis(2-ethylhexyl)phthalate	50.9	J	2.3	5.0	"	50.0	6.0	89.8	39-138	25.1	33
Di-n-octyl phthalate	48.4		2.7	5.0	"	50.0	U	96.8	28-158	22.5	46
Benzo (b) fluoranthene	105		1.1	5.0	"	50.0	U	210	49-155	79.8	22
Benzo (k) fluoranthene	89.4		1.1	5.0	"	50.0	U	179	43-149	76.8	18
Benzo (a) pyrene	36.4		0.73	5.0	"	50.0	U	72.8	31-160	13.1	47
Indeno (1,2,3-cd) pyrene	41.2		0.30	5.0	"	50.0	U	82.4	52-171	5.74	37
Dibenz (a,h) anthracene	40.8		0.63	5.0	"	50.0	U	81.6	12-226	4.77	41
Benzo (g,h,i) perylene	39.3		0.93	5.0	"	50.0	U	78.6	25-204	6.03	37
Surrogate: 2-Fluorophenol	17.5				"	50.0		35.0	18-128		
Surrogate: Phenol-d6	28.6				"	50.0		57.2	12-133		
Surrogate: Nitrobenzene-d5	16.9				"	50.0		33.8	5-158		
Surrogate: 2-Fluorobiphenyl	32.0				"	50.0		64.0	29-128		
Surrogate: 2,4,6-Tribromophenol	42.7				"	50.0		85.4	32-148		
Surrogate: Terphenyl-d14	47.1				"	50.0		94.2	62-140		





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Project:Himco Dump  
Project Number:2003SY01  
Project Manager:Howard Pham

**Reported:**  
Dec-05-02 14:50

### Notes and Definitions

- J The identification of the analyte is acceptable; the reported value is an estimate.
- R The presence or absence of the analyte cannot be determined from the data due to severe quality control problems. The data are rejected and considered unuseable.
- UJ The analyte was not detected at or above the reported value. The reported value is an estimate.
- U Not Detected
- NR Not Reported

Sample E2K0101-06 was used as the MS1/MSD1 samples. The MS1 and MSD1 samples were each spiked with 500ul of the 100ug/ml ABN LCS standard solution.

In the MS1 sample all the percent recoveries were within limits except Pentachlorophenol (188%) and 3,3-Dichlorobenzidine(26%).

The analyte 2,4-Dinitrophenol was not detected in either MS1 or MSD1. This analyte was flagged "Q" for detected and "R" for non-detected compounds in the native sample.

In the MSD1 sample all the percent recoveries were within limits except:

bis(2-Chloroethyl)ether (10%), 1,3-Dichlorobenzene(11%), 1,4-Dichlorobenzene(12%), 1,2-Dichlorobenzene(16%), bis(2-chloroisopropyl)ether(24%), 2-Nitrophenol(38%) N-Nitroso-di-n-propylamine(32%), Nitrobenzene(32%), Isophorone(42%), bis(2-Chloroethoxy)methane(44%), 1,2,4-Trichlorobenzene(32%) 4-Chloroaniline (26%), Pentachlorophenol(162%) and 3,3-Dichlorobenzidine (18%). All these compounds percent recoveries were within limits in the MS1 except the analytes Pentachlorophenol and 3,3-Dichlorobenzidine were outside limits in both MS1 and MSD1. These compounds were flagged "J" for detected and "UJ" for non-detected compounds in the native sample.

The relative percent differences(RPD) were within limits except bis(2-Chloroethyl)ether (117%), 1,3-Dichlorobenzene(123%), 1,4-Dichlorobenzene(120%), 1,2-Dichlorobenzene (103%), bis(2-chloroisopropyl)ether(83%), Hexachloroethane(97%), N-Nitroso-di-n-propylamine(64%), Nitrobenzene(58%), Isophorone(42%), 2,4-Nitrophenol (51%),4-Chloroaniline (76%) and bis(2-Ethylhexyl)phthalate(29%). No other problems were observed.

#### **IV.SAMPLES RESULTS:**

In the sample E2k0101-02 contained bis(2-Ethylhexyl)phthalate(5.0ug/L)

In the sample E2k0101-04 contained bis(2-Ethylhexyl)phthalate(7.0ug/L)

In the sample E2k0101-05 contained bis(2-Ethylhexyl)phthalate(4.0ug/L)

In the sample E2k0101-06 contained bis(2-Ethylhexyl)phthalate(6.0ug/L)

The raw /processed data files, method files and sequence files are archived on R5CRL\VOL2\XNGUYEN\2002.

Tentatively Identified Compounds (TICs) have been listed for each sample. These compounds listed on the Enviroforms were selected based on the best Mass spectrum match of the unknown compound to compounds in the library. These compounds are suggested and not confirmed. LIMS reports will be generated as soon as possible.

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